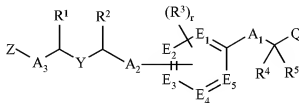


Amendments to the Claims

1. (Currently Amended). A compound having a formula I,



I

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ is: ~~a bond, CH₂, O or S, and wherein A₁ and R⁴ or A₁ and R⁵ together being a 3- to 6-membered carbocycl~~ when A₁ is a carbon;

A₂ and A₃ are independently: CH₂, O or S;

E₁, E₂, E₃, E₄ and E₅ are each CH or substituted carbon bearing A₂ and R³; or at least one of E₁, E₂, E₃, E₄ and E₅ is nitrogen and each of others being CH or substituted carbon bearing A₂ and R³;

Q is: -C(O)OR⁶ ~~or R^{6A}~~;

Y is: ~~a bond~~ or C₁-C₆ alkyl or C₃-C₆ cycloalkyl;

Z is: a) phenyl-T-pyridinyl or naphthyl-T-pyridinyl each optionally substituted with one or more R⁷; wherein T is a single bond, C or O; aryl;

b) ~~a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S;~~

c) ~~bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or~~

d) ~~bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R⁷;~~

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: ~~1, 2, 3, or 4;~~

R¹ and R² are each independently:

hydrogen,

haloalkyl,

C₁-C₆ alkyl,

(CH₂)_nC₃-C₈ cycloalkyl, or

~~R¹ and R² form a 4 to 8 membered nonaromatic carbocyclic ring; and~~

~~wherein at least one of R¹ and R² is alkyl or cycloalkyl; and;~~

R³ is: hydrogen,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

C₁-C₆ alkyl,

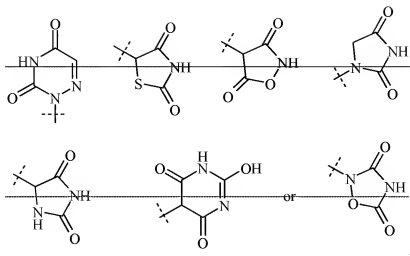
C₁-C₆ alkoxy, or

C₃-C₈ cycloalkyl;

~~R⁴ and R⁵ are each independently: hydrogen or C₁-C₆ alkyl;~~

R⁶ is: hydrogen, C₁-C₆ alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,



R⁷ is: hydrogen,

oxo,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

arylalkyl,

aminoalkyl,

C₁-C₆ alkyl,

C₁-C₆ alkoxy,

(CH₂)_nC₃-C₈ cycloalkyl,

C(O)R⁹,

C(O)OR⁹,

C(=NOR⁸)R⁹,

CR⁸(OH)R⁹,

C[=C(R⁸)₂]R⁹,

OR⁹,

SR⁹ or

S(O)_pR⁹;

R⁸ is: hydrogen or C₁-C₆ alkyl; and

R⁹ is: hydrogen,

C₁-C₆ alkyl,

C₃-C₈ cycloalkyl,

aryl,

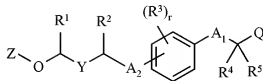
heteroaryl or

heterocyclyl,

wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C₁-C₆ alkyl, C₁-C₆ alkoxy and C₃-C₈ cycloalkyl.

2. (Currently Amended). The compound of Claim 1, wherein the compound is represented by a compound of formula II,



II

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ is: ~~a bond, CH₂, O or S; and wherein A₁ and R⁴ or A₁ and R⁵ together being a 3- to 6-membered carbocyclyl when A₁ is a carbon;~~

A₂ is: O or S or CH₂;

Q is: -C(O)OR⁶, ~~or R^{6a};~~

Y is: a bond or; C₁-C₆ alkyl or C₃-C₈ cycloalkyl;

Z is: phenyl-T-pyridinyl or naphthyl-T-pyridinyl each optionally substituted with one or more R⁷; wherein T is a single bond, C or O; ~~a~~ aryl;

b) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S;

- e) ~~bi-aryl~~, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or ~~—~~
- d) ~~bi-heteroaryl~~, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R^2 ;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: ~~1, 2, 3, or 4;~~

R^1 and R^2 are each independently:

hydrogen,

~~haloalkyl,~~

C_1 - C_6 alkyl;

~~$(CH_2)_n$, C_3 - C_8 cycloalkyl, or~~

R^1 and R^2 form a 4 to 8 membered nonaromatic carbocyclic ring; and

wherein at least one of R^1 and R^2 is alkyl or cycloalkyl; and;

R^3 is: hydrogen,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

C_1 - C_6 alkyl,

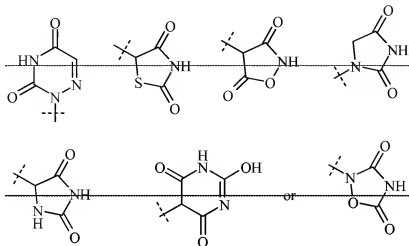
C_1 - C_6 alkoxy, or

C_3 - C_8 cycloalkyl;

R^4 and R^5 are each independently: hydrogen or C_1 - C_6 alkyl;

R^6 is: hydrogen, C_1 - C_6 alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole:



R^7 is: hydrogen,

oxo,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

aryloxy,

arylalkyl,

aminoalkyl,

C_1 - C_6 alkyl,

C_1 - C_6 alkoxy,

$(CH_2)_n$ - C_3 - C_8 cycloalkyl,

$C(O)R^9$,

$C(O)OR^9$,

$C(=NOR^8)R^9$,

$CR^8(OH)R^9$,

$C[=C(R^8)_2]R^9$,

OR^9 ,

SR^9 or

$S(O)_pR^9$;

R⁸ is: hydrogen or C₁-C₆ alkyl; and

R⁹ is: hydrogen,

C₁-C₆ alkyl,

C₃-C₈ cycloalkyl,

aryl,

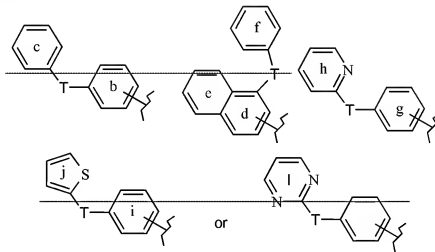
heteroaryl or

heterocyclyl,

wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C₁-C₆ alkyl, C₁-C₆ alkoxy and C₃-C₈ cycloalkyl.

3. (Currently Amended). The compound of Claim 2, wherein Z is an optionally substituted phenyl or naphthyl, furanyl, imidazolyl, indolyl, oxazolyl, isoxazolyl, pyridyl, pyrrolyl, thiazolyl, thiophenyl, benzofuranyl, benzothiophenyl, benzoisoxazolyl, quinolinyl, isoquinolinyl or a structural formula selected from following:



wherein T is:

a bond, $-(CH_2)_qO-$, $-O(CH_2)_q-$, $-C(O)(CH_2)_q-$, $-(CH_2)_qC(O)-$, $-(CH_2)_qS-$, $-S(CH_2)_q-$, $-S(O)_2-$, $-(C_1-C_3\text{ alkyl})-$, $-(CH_2)_qC(=CH_2)-$, $-C(=CH_2)(CH_2)_q-$, $-(CH_2)_qC(=NOH)-$, $-C(=NOH)(CH_2)_q-$, $-(CH_2)_qC(=NOCH_3)-$, $-C(=NOCH_3)(CH_2)_q-$, $-CH(OH)(CH_2)_q-$, or $-(CH_2)_qCH(OH)-$,

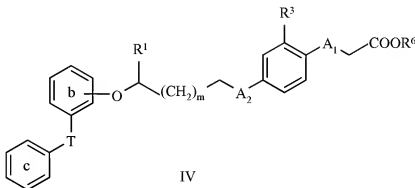
q is: 0, 1, 2 or 3; and

rings b to h are each optionally substituted with one or more groups independently selected from the group consisting of:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1 - C_6 alkyl, C_1 - C_6 alkoxy and $(CH_2)_hC_3$ - C_8 cycloalkyl.

4. (Canceled)

5. (Withdrawn). The compound of Claim 2, wherein the compound is represented by structural formula IV,



or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A_1 and A_2 are respectively:

- O and O,
- CH_2 and O,
- CH_2 and S,
- O and S or
- S and O;

m is: 1 or 2;

R^1 is: C_1 - C_3 alkyl;

R^3 is: hydrogen, halo or C_1 - C_6 alkyl;

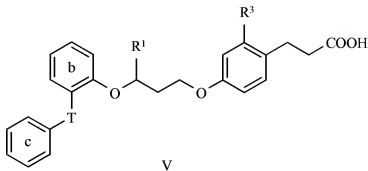
R^6 and R^9 are each independently: hydrogen or C_1 - C_6 alkyl;

T is: a bond, -O-, -C(O)-, -S(O)-S(O)₂-, -C(=CH₂)-, -C(=NOH)- or -CH(OH)-; and

rings b and c are each optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, $S(O)_2R^9$, C_1 - C_6 alkyl, C_1 - C_6 alkoxy and $(CH_2)_hC_3$ - C_8 cycloalkyl.

6. (Withdrawn). The compound of Claim 5, wherein the compound is represented by structural formula V,



or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

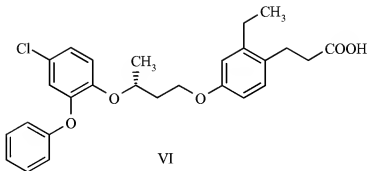
T is: a bond, -O- or -C(O)-;

R¹ is: methyl, ethyl or cyclopropyl;

R³ is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, methyl, ethyl, isopropyl, N(CH₃)₂, S(O)₂CH₃, methoxy and cyclopropyl.

7. (Withdrawn). The compound of Claim 6, wherein the compound is represented by a structural formula VI,

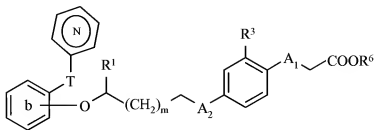


or a pharmaceutically acceptable salt, solvate or hydrate thereof.

8. (Canceled)

9. (Canceled)

10. (Currently amended). The compound of Claim 2, wherein the compound is represented by structural formula VIII,



VIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ and A₂ are respectively:

- O and O,
- CH₂ and O,
- CH₂ and S,
- O and S or
- S and O;

m is: 1 or 2;

R¹ is: C₁-C₃ alkyl; and

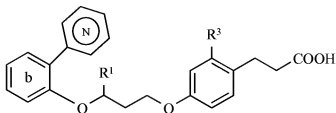
R³ is: hydrogen, halo or C₁-C₆ alkyl;

R⁶ and R⁹ are each independently: hydrogen or C₁-C₆ alkyl;

T is: a ~~bond or bond~~, -O-, -C(O)-, -S(O)-, -S(O)₂-, -C(=CH₂)-, -C(=NOH)- or -CH(OH)-; and ring b is optionally substituted with one or more groups independently selected from:

- hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, S(O)₂R⁹, C₁-C₆ alkyl, C₁-C₆ alkoxy and (CH₂)_nC₃-C₈ cycloalkyl.

11. (Previously Presented). The compound of Claim 10, wherein the compound is represented by structural formula IX,



IX

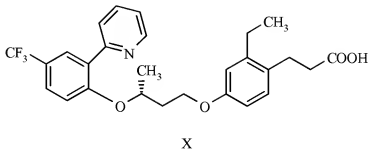
or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

R¹ is C₁-C₃ alkyl;

R³ is: hydrogen, halo or C₁-C₄ alkyl;

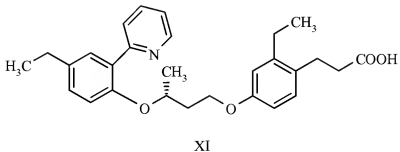
ring b is optionally substituted with one or more groups independently selected from the group consisting of: hydrogen, halo, haloalkyl, haloalkyloxy and C₁-C₆ alkyl.

12. (Previously Presented). The compound of Claim 11, wherein the compound is represented by structural formula X,

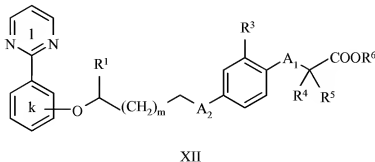


or a pharmaceutically acceptable salt, solvate or hydrate thereof.

13. (Previously Presented). The compound of Claim 11, wherein the compound is represented by structural formula XI,



14. (Withdrawn). The compound of Claim 2, wherein the compound is represented by structural formula XII,



or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ and A₂ are respectively:

- O and O,
- CH₂ and O,
- CH₂ and S,
- O and S or
- S and O;

m is: 1 or 2;

R¹ is: C₁-C₃ alkyl; and

R³ is: hydrogen, halo or C₁-C₆ alkyl;

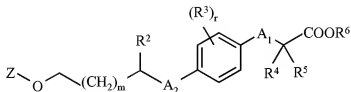
R⁴, R⁵, R⁶ and R⁹ are each independently: hydrogen or C₁-C₆ alkyl;

rings k and l are each optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, S(O)₂R⁹, C₁-C₆ alkyl, C₁-C₆ alkoxy and (CH₂)_mC₃-C₈ cycloalkyl.

15. (Canceled)

16. (Previously Presented). The compound of Claim 2, wherein the compound is represented by structural formula XIII,



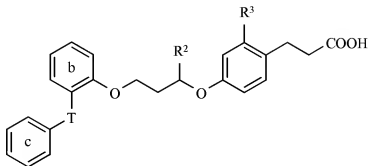
XIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein

m is 1, 2, 3, or 4.

17. (Canceled).

18. (Withdrawn). The compound of Claim 16, wherein the compound is represented by structural formula XV,



XV

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

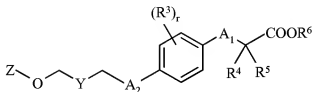
T is: a bond, O or C(O);

R² is: methyl, ethyl or cyclopropyl;

R³ is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, N(CH₃)₂, S(O)₂CH₃, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

19. (Currently Amended). The compound of Claim 2, wherein the compound is represented by structural formula XVI,

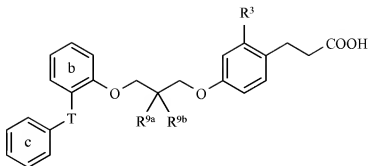


XVI

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein Y is a branched alkyl or C₃-C₆-cycloalkyl.

20. (Canceled).

21. (Withdrawn). The compound of Claim 19, wherein the compound structural formula XVIII,



XVIII

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

T is: a bond, O or C(O);

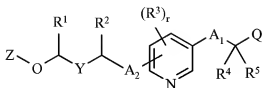
R³ is: methyl or ethyl;

R^{9a} and R^{9b} are each independently hydrogen, methyl or ethyl, wherein at least one of R^{9a} and R^{9b} being methyl or ethyl;

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, S(O)₂CH₃, N(CH₃)₂, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

22. (Canceled).

23. (Withdrawn). The compound of Claim 1, wherein the compound is a compound of formula XX,



XX

or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ is: a bond, CH₃, O or S, and wherein A₁ and R⁴ or A₁ and R⁵ together being a 3- to 6-membered carbocyclyl when A₁ is a carbon;

A₂ is: O or S or CH₂;

Q is: -C(O)OR⁶, or R^{6A};

Y is: a bond, C₁-C₆ alkyl or C₃-C₆ cycloalkyl;

- Z is:
- a) aryl;
 - b) a 5- to 10-membered heteroaryl wherein the heteroaryl containing at least one heteroatom selected from N, O or S,
 - c) bi-aryl, wherein biaryl being defined as aryl substituted with another aryl or aryl substituted with heteroaryl, or
 - d) bi-heteroaryl, wherein bi-heteroaryl being defined as heteroaryl substituted with another heteroaryl, or heteroaryl substituted with aryl, and wherein aryl, heteroaryl, bi-aryl and bi-heteroaryl being optionally substituted with one or more groups independently selected from R⁷;

n is: 1, 2, 3, 4, 5 or 6

p is: 1 or 2;

r is: 1, 2, 3, or 4;

R¹ and R² are each independently:

hydrogen,

haloalkyl,

C₁-C₆ alkyl,

(CH₂)_nC₃-C₈ cycloalkyl, or

R¹ and R² form a 4- to 8-membered nonaromatic carbocyclic ring; and wherein at least one of R¹ and R² is alkyl or cycloalkyl, and;

R³ is: hydrogen,

nitro,

cyano,

hydroxyl,

halo,

haloalkyl,

haloalkyloxy,

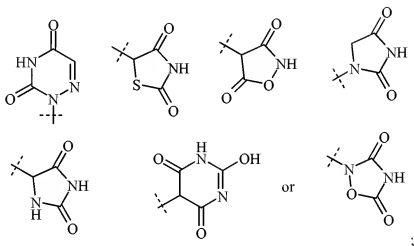
aryloxy,

C₁-C₆ alkyl,
 C₁-C₆ alkoxy, or
 C₃-C₈ cycloalkyl;

R⁴ and R⁵ are each independently: hydrogen or C₁-C₆ alkyl;

R⁶ is: hydrogen, C₁-C₆ alkyl or aminoalkyl;

R^{6A} is: carboxamide, sulfonamide, acylsulfonamide, tetrazole,



R⁷ is: hydrogen,

oxo,
 nitro,
 cyano,
 hydroxyl,
 halo,
 haloalkyl,
 haloalkyloxy,
 aryloxy,
 arylalkyl,
 aminoalkyl,
 C₁-C₆ alkyl,
 C₁-C₆ alkoxy,
 (CH₂)_nC₃-C₈ cycloalkyl,
 C(O)R⁹,

$C(O)OR^9$,
 $C(=NOR^8)R^9$,
 $CR^8(OH)R^9$,
 $C[=C(R^8)_2]R^9$,
 OR^9 ,
 SR^9 or
 $S(O)_pR^9$;

R^8 is: hydrogen or C_1 - C_6 alkyl; and

R^9 is: hydrogen,

C_1 - C_6 alkyl,

C_3 - C_8 cycloalkyl,

aryl,

heteroaryl or

heterocyclyl,

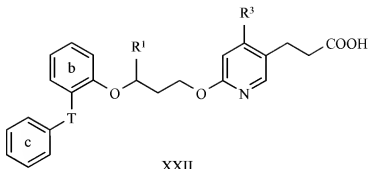
wherein alkyl, cycloalkyl, aryl, heteroaryl or heterocyclyl being optionally substituted with one or more substituents selected from the group consisting of:

hydrogen, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, oxo, C_1 - C_6 alkyl, C_1 - C_6 alkoxy and C_3 - C_8 cycloalkyl.

24. (Canceled).

25. (Canceled).

26. (Withdrawn). The compound of Claim 23, wherein the compound is a compound of structural formula XXII,



or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

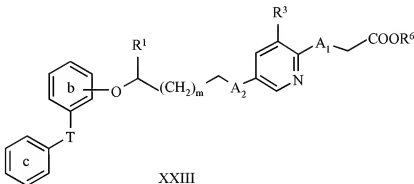
T is: a bond, -O- or -C(O)-;

R¹ is: methyl, ethyl or cyclopropyl;

R³ is: methyl or ethyl; and

rings b and c are each optionally substituted with one or more substituent independently selected from the group consisting of: hydrogen, Cl, Br, CF₃, OCF₃, S(O)₂CH₃, N(CH₃)₂, methyl, ethyl, isopropyl, methoxy and cyclopropyl.

27. (Withdrawn). The compound of Claim 1, wherein the compound is a compound of structural formula XXIII,



or a pharmaceutically acceptable salt, solvate, hydrate or stereoisomer thereof, wherein:

A₁ and A₂ are respectively:

O and O,

CH₂ and O,

CH₂ and S,

O and S or

S and O;

m is: 1, 2, 3 or 4;

R¹ is: C₁-C₃ alkyl; and

R³ is: hydrogen, halo or C₁-C₆ alkyl;

R⁶ and R⁹ are each independently: hydrogen or C₁-C₆ alkyl;

T is: a bond, -O-, -C(O)-, -S(O)-S(O)₂-, -C(=CH₂)-, -C(=NOH)- or -CH(OH)-; and

rings b and c are each optionally substituted with one or more groups independently selected from:

hydrogen, oxo, nitro, cyano, hydroxyl, halo, haloalkyl, haloalkyloxy, aryloxy, arylalkyl, aminoalkyl, S(O)₂R⁹, C₁-C₆ alkyl, C₁-C₆ alkoxy and (CH₂)_nC₃-C₈ cycloalkyl.

28. (Canceled).

29. (Currently Amended). A compound of Claim 1 selected from the group consisting of:

	Structure	Name
1		3-[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
2		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-acetic acid
3		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

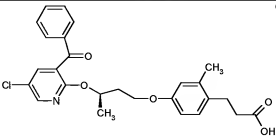
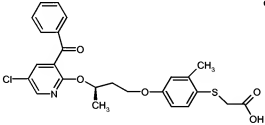
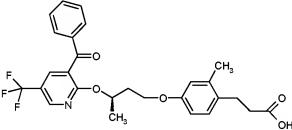
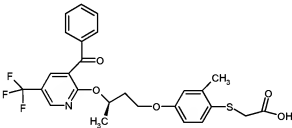
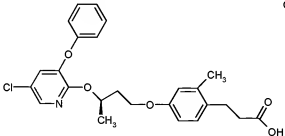
	Structure	Name
4		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
5		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenoxy}-acetic acid
6		3-[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenyl]-propionic acid
7		2-[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy]-2-methyl-propionic acid
8		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-acetic acid

	Structure	Name
9		3-[4-{3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
10		3-[4-{3-(2-Benzoyl-4-cyclopropyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
11		3-[4-{3-(2-Benzoyl-4-trifluoromethyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
12		3-[4-{3-(2-Benzoyl-4-chloro-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
13		3-[4-{3-(2-Benzoyl-4-chloro-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid

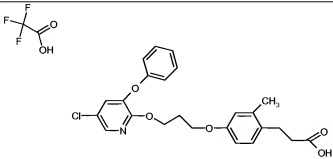
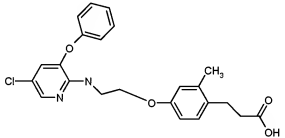
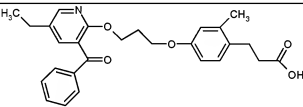
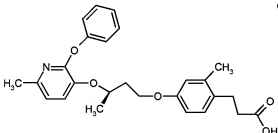
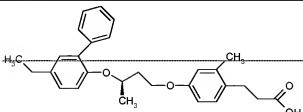
	Structure	Name
14	<p>Chiral</p>	3-[4-{3-(2-Benzoyl-4-methoxy-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
15	<p>Chiral</p>	3-[4-{3-(2-Benzoyl-4-fluoro-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
16	<p>Chiral</p>	3-[4-{3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
17		4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenylsulfanyl]-acetic acid
18		3-(4-[3-[4-Ethyl-2-(hydroxy-phenyl-methyl)-phenoxy]-butoxy]-2-methyl-phenyl)-propionic acid

	Structure	Name
19		3-(4-{3-[4-Ethyl-2-(hydroxyimino-phenyl)-methyl]-phenoxy}-butoxy)-2-methyl-phenyl)-propionic acid
20		3-(4-{3-[4-Ethyl-2-(methoxyimino-phenyl)-methyl]-phenoxy}-butoxy)-2-methyl-phenyl)-propionic acid
21		3-{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl)-propionic acid
22		{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
23		3-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-2-methyl-phenyl)-propionic acid
24		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl)-propionic acid

	Structure	Name
25		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
26		3-{4-[3-(2-Cyclopentanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
27		2-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
28		2-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
29		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
30		{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

	Structure	Name
31		3-{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
32		{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
33		3-{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
34		{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
35		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid

	Structure	Name
36		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-ethyl-phenyl}-propionic acid
37		{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
38		3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
39		3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
40		3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
41		3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-propoxy]-phenyl}-propionic acid (trifluoroacetic acid salt)

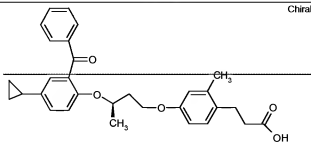
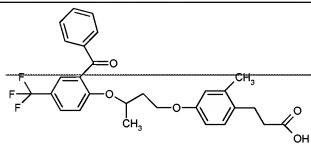
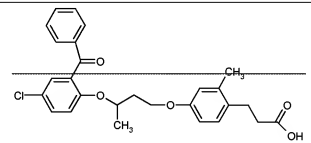
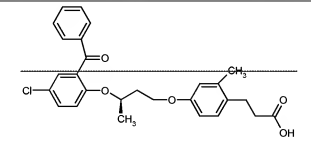
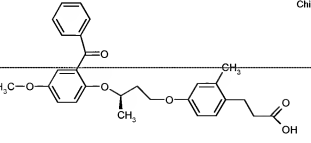
	Structure	Name
42		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
43		3-{4-[2-(5-Chloro-3-phenoxy-pyridin-2-ylamino)-ethoxy]-2-methyl-phenyl}-propionic acid
44		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
45		Chiral 3-{2-Methyl-4-[3-(6-methyl-2-phenoxy-pyridin-3-yloxy)-butoxy]-phenyl}-propionic acid
46		3-{4-[3-(5-Ethyl biphenyl-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid

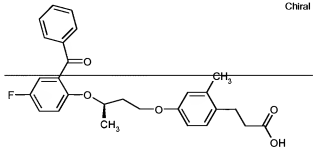
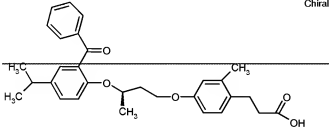
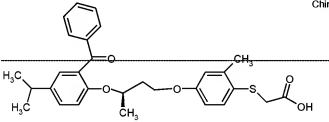
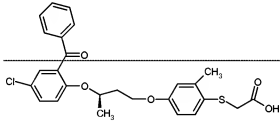
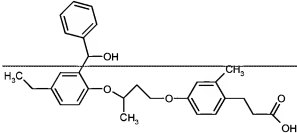
	Structure	Name
47		3-{4-[3-(4-Ethyl-2-oxazol-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
48		3-{4-[3-(4-Ethyl-2-thiazol-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
49		3-{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
50		{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
51		3-{2-Ethyl-4-[3-(4-ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-phenyl}-propionic acid
52		3-{4-[3-(4-Chloro-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

	Structure	Name
53		3-{2-Methyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
54		3-{2-Ethyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
55		3-{4-[3-(4-Ethyl-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
56		3-{4-[3-(4-Chloro-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
57		3-{4-[3-(4-Ethyl-2-pyridin-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
58		3-{2-Methyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid

	Structure	Name
59		3-{[2-Ethyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl]-propionic acid
60		3-{[4-[3-(2-Benzof[d]isoxazol-3-yl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
61		3-{[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
62		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy}-acetic acid
63		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
64		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

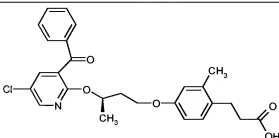
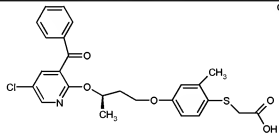
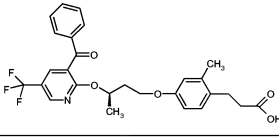
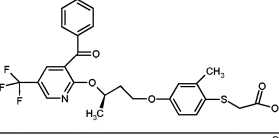
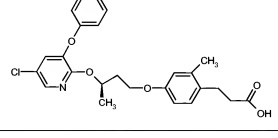
	Structure	Name
65		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenoxy}-acetic acid
66		3-[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butylsulfanyl]-2-methyl-phenyl]-propionic acid
67		2-[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenoxy]-2-methyl-propionic-acid
68		{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-acetic acid
69		3-[4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic-acid

	Structure	Name
70		3-[4-{3-(2-Benzoyl-4-cyclopropyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
71		3-[4-{3-(2-Benzoyl-4-trifluoromethyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
72		3-[4-{3-(2-Benzoyl-4-chloro-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
73		3-[4-{3-(2-Benzoyl-4-chloro-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
74		3-[4-{3-(2-Benzoyl-4-methoxy-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid

	Structure	Name
75		3-[4-{3-(2-Benzoyl-4-fluoro-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
76		3-[4-{3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy}-2-methyl-phenyl]-propionic acid
77		4-[3-(2-Benzoyl-4-isopropyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl]-acetic acid
78		4-[3-(2-Benzoyl-4-chloro-phenoxy)-butoxy]-2-methyl-phenylsulfanyl]-acetic acid
79		3-(4-{3-[4-Ethyl-2-(hydroxy-phenyl-methyl)-phenoxy]-butoxy}-2-methyl-phenyl)-propionic acid

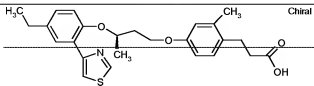
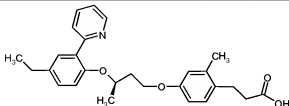
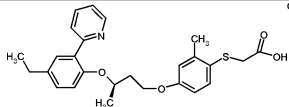
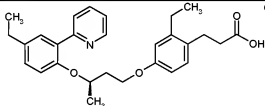
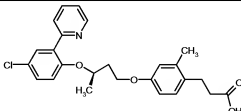
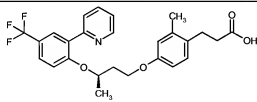
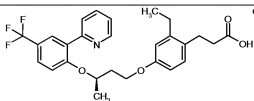
	Structure	Name
80		3-(4-{3-[4-Ethyl-2-(hydroxyimino-phenyl)-methyl]-phenoxy}-butoxy)-2-methyl-phenyl)-propionic acid
81		3-(4-{3-[4-Ethyl-2-(methoxyimino-phenyl)-methyl]-phenoxy}-butoxy)-2-methyl-phenyl)-propionic acid
82		3-{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
83		{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-phenylsulfanyl}-acetic acid
84		3-{4-[3-(4-Ethyl-2-isobutyryl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
85		3-{4-[3-(2-Cyclopropanecarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

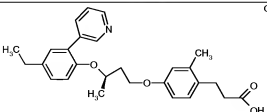
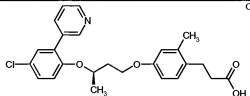
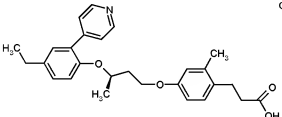
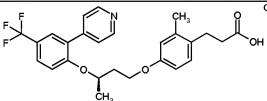
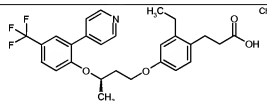
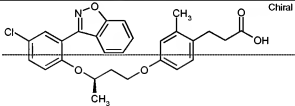
	Structure	Name
86		3-{4-[3-(2-Cyclopropylcarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
87		3-{4-[3-(2-Cyclopentylcarbonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
88		2-{4-[3-(4-Ethyl-2-isobutyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
89		2-{4-[3-(2-Cyclopropylcarbonyl-4-ethyl-phenoxy)-butoxy]-phenoxy}-2-methyl-propionic acid
90		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
91		{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid

	Structure	Name
92		3-{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
93		{4-[3-(3-Benzoyl-5-chloro-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
94		3-{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
95		{4-[3-(3-Benzoyl-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
96		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid

	Structure	Name
97		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-ethyl-phenyl}-propionic acid
98		{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
99		3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
100		3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
101		3-{2-Ethyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-butoxy]-phenyl}-propionic acid
102		3-{2-Methyl-4-[3-(3-phenoxy-5-trifluoromethyl-pyridin-2-yloxy)-propoxy]-phenyl}-propionic acid (trifluoroacetic acid salt)

	Structure	Name
103		3-{4-[3-(5-Chloro-3-phenoxy-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
104		3-{4-[2-(5-Chloro-3-phenoxy-pyridin-2-ylamino)-ethoxy]-2-methyl-phenyl}-propionic acid
105		3-{4-[3-(3-Benzoyl-5-ethyl-pyridin-2-yloxy)-propoxy]-2-methyl-phenyl}-propionic acid
106		3-{2-Methyl-4-[3-(6-methyl-2-phenoxy-pyridin-3-yloxy)-butoxy]-phenyl}-propionic acid
107		3-{4-[3-(5-Ethyl-biphenyl-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic acid
108		3-{4-[3-(4-Ethyl-2-oxazol-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

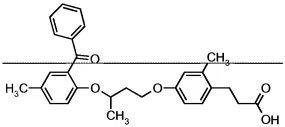
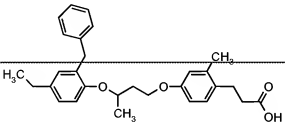
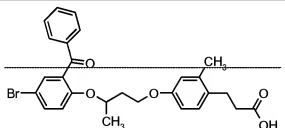
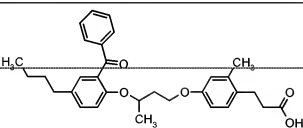
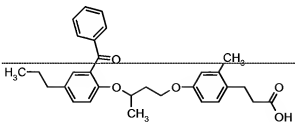
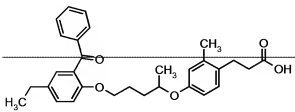
	Structure	Name
109		3-{4-[3-(4-Ethyl-2-thiazol-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
110		3-{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
111		{4-[3-(4-Ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
112		3-{2-Ethyl-4-[3-(4-ethyl-2-pyridin-2-yl-phenoxy)-butoxy]-phenyl}-propionic acid
113		3-{4-[3-(4-Chloro-2-pyridin-2-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
114		3-{2-Methyl-4-[3-(2-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
115		3-{2-Ethyl-4-[3-(2-pyridin-2-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid

	Structure	Name
116	 <p>Chiral</p>	3-{4-[3-(4-Ethyl-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
117	 <p>Chiral</p>	3-{4-[3-(4-Chloro-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
118	 <p>Chiral</p>	3-{4-[3-(4-Ethyl-2-pyridin-4-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
119	 <p>Chiral</p>	3-{2-Methyl-4-[3-(2-trifluoromethyl-4-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
120	 <p>Chiral</p>	3-{2-Ethyl-4-[3-(2-pyridin-4-yl-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
121	 <p>Chiral</p>	3-{4-[3-(2-chloro-phenoxyl)-butoxy]-2-methyl-phenyl}-propionic acid

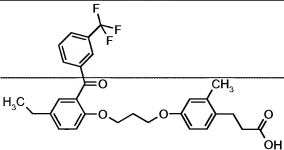
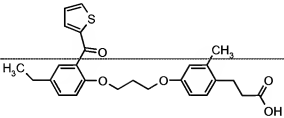
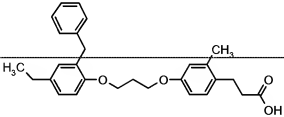
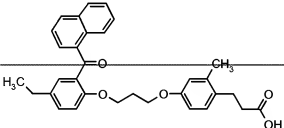
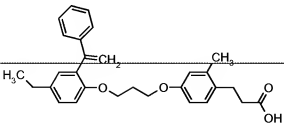
	Structure	Name
122		(R)-[4-{3-(4-ethyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
123		(R)-[4-{3-(2-benzoyl-4-methyl phenoxy) butoxy]-2-methyl phenylsulfanyl}-acetic acid
124		(R)-[4-{3-(2-benzoyl-4-trifluoromethoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
125		{4-[3-(2-benzoyl-4-ethyl-phenoxy)-hexyloxy]-2-methyl-phenylsulfanyl}-acetic acid
126		3-[4-[3-(2-benzoyl-4-ethyl-phenoxy)-hexyloxy]-2-methyl-phenyl]-propionic acid

	Structure	Name
127		(R)-3-{4-[3-(4-ethyl-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic-acid
128		(R)-3-{4-[3-[4-ethyl-2-(1-phenyl-vinyl)-phenoxy]-butoxy]-2-methyl-phenyl}-propionic-acid
129		(R)-3-{4-[3-[4-ethyl-2-(1-methyl-1-phenyl-ethyl)-phenoxy]-butoxy]-2-methyl-phenyl}-propionic acid
130		(R)-3-{4-[3-(2-benzoyl-4-methyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
140		(R)-3-{4-[3-[4-ethyl-2-(1-phenyl-ethyl)-phenoxy]-butoxy]-2-methyl-phenyl}-propionic-acid
141		(R)-3-{4-[3-[4-ethyl-2-(pyridine-2-carbonyl)-phenoxy]-butoxy]-2-methyl-phenyl}-propionic acid

	Structure	Name
142		3-(2-methyl-4-((3-(2-(trifluoromethoxy)phenyl)propionyloxy)phenyl)butoxy)-(thiophene-2-carbonyl)-4-trifluoromethoxy-phenyl propionic acid
143		3-(4-((3-(4-ethyl-2-(thiophene-2-carbonyl)phenoxy)butoxy)-2-methylphenyl)propionic acid
144		3-(4-((3-(4-ethyl-2-(naphthalene-1-carbonyl)phenoxy)butoxy)-2-methylphenyl)propionic acid
145		3-(4-((3-(4-ethyl-2-(1-phenylvinyl)phenoxy)butoxy)-2-methylphenyl)propionic acid
146		3-(4-((3-(2-benzoylphenoxy)butoxy)-2-methylphenyl)propionic acid

	Structure	Name
147		3-[4-[3-(2-benzoyl-4-methyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
148		3-[4-[3-(2-benzyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
149		3-[4-[3-(2-benzoyl-4-bromo-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
150		3-[4-[3-(2-benzoyl-4-butyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
151		3-[4-[3-(2-benzoyl-4-propyl-phenoxy)-butoxy]-2-methyl-phenyl]-propionic acid
152		3-[4-[4-(2-benzoyl-4-ethyl-phenoxy)-1-methyl-butoxy]-2-methyl-phenyl]-propionic acid

	Structure	Name
153		3-[4-[4-(2-benzoyl-4-ethyl-phenoxy)-pentyloxy]-2-methyl-phenyl]-propionic acid
154		3-[4-[3-(2-benzoyl-4-ethyl-phenoxy)-2-methyl-propoxy]-2-methyl-phenyl]-propionic acid
155		3-[4-[3-(2-benzoyl-4-ethyl-phenoxy) propoxy]-2-methyl phenyl]-propionic acid
156		3-[4-[3-[4-ethyl-2-(4-fluoro-benzoyl)-phenoxy]-propoxy]-2-methyl-phenyl]-propionic acid
157		3-[4-[3-[4-ethyl-2-(2-trifluoromethyl-benzoyl)-phenoxy]-propoxy]-2-methyl phenyl]-propionic acid

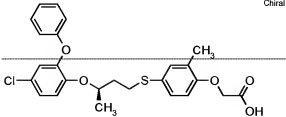
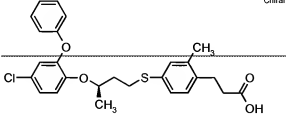
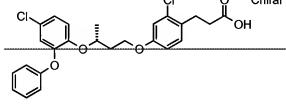
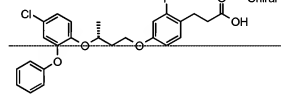
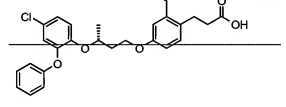
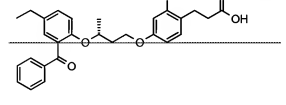
	Structure	Name
158		3-(4-{3-[4-ethyl-2-(3-trifluoromethyl-benzoyl)-phenoxy]propoxy}-2-methyl-phenyl)-propionic acid
159		3-(4-{3-[4-ethyl-2-(thiophene-2-carbonyl)-phenoxy]propoxy}-2-methyl-phenyl)-propionic acid
160		3-(4-{3-[2-benzyl-4-ethyl-phenoxy]propoxy}-2-methyl-phenyl)-propionic acid
161		3-(4-{3-[4-ethyl-2-(naphthalene-1-carbonyl)-phenoxy]propoxy}-2-methyl-phenyl)-propionic acid
162		3-(4-{3-[4-ethyl-2-(1-phenyl-vinyl)-phenoxy]propoxy}-2-methyl-phenyl)-propionic acid

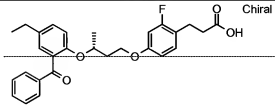
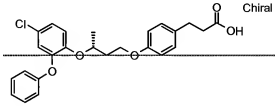
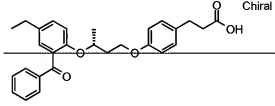
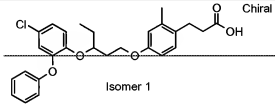
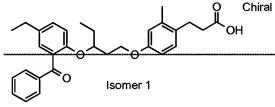
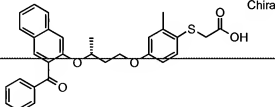
	Structure	Name
163		2-[4-[3-(2-benzoyl-4-ethyl-phenoxy)-butoxy]-phenoxyl-2-methyl-propionic-acid
164		2-[4-[3-(2-benzoyl-4-ethyl-phenoxyl-2-methyl-propoxy)-phenoxyl-2-methyl-propionic acid
165		2-[4-[3-(2-benzyl-4-ethyl-phenoxyl)-butoxy]-phenoxyl-2-methyl-propionic-acid
166		2-[4-[3-(2-benzoyl-4-bromo-phenoxyl)-butoxy]-phenoxyl-2-methyl-propionic-acid

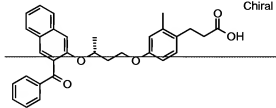
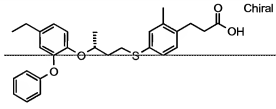
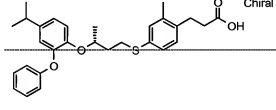
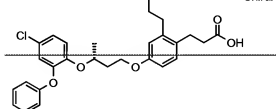
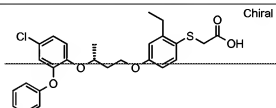
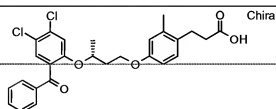
	Structure	Name
167		2-[4-{3-(2-benzoyl-4-butyl-phenoxy)-butoxy}-phenoxy]-2-methyl-propionic-acid
168	<p>Chiral</p>	(R)-3-[4-{3-(4-chloro-2-phenoxy-phenoxy)-butoxy}-2-methyl-phenyl]-propionic-acid
169	<p>Chiral</p>	(R)-3-[2-methyl-4-{3-(2-phenoxy-4-trifluoromethyl-phenoxy)-butoxy}-phenyl]-propionic-acid
170	<p>Chiral</p>	(R)-3-[2-methyl-4-{3-(2-phenoxy-4-trifluoromethoxy-phenoxy)-butoxy}-phenyl]-propionic-acid
171	<p>Chiral</p>	(R)-3-[2-methyl-4-{3-(4-methyl-2-phenoxy-phenoxy)-butoxy}-phenyl]-propionic-acid

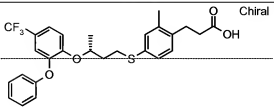
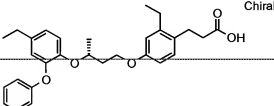
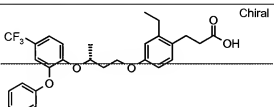
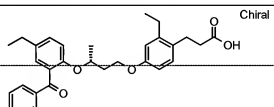
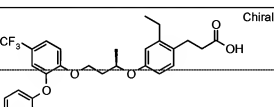
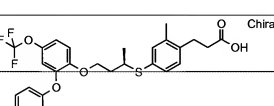
	Structure	Name
172		(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-2-methyl-phenylsulfanyl}-acetic acid
173		3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-propoxy]-2-methyl-phenyl}-propionic acid
174		(R)-3-{4-[3-(2-benzothiophen-3-yl-4-chloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
175		(R)-3-{4-[3-(4-chloro-2-pyridin-3-yl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
176		(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-2,2-difluoro-propionic acid
177		%)(R)-3-{3-bromo-4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid

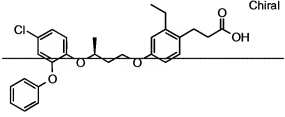
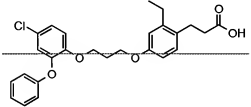
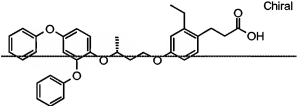
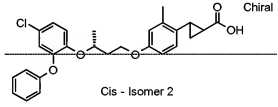
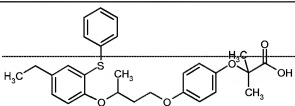
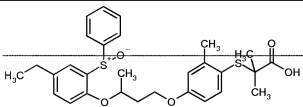
	Structure	Name
178	<p>Chiral</p>	(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-3-methyl-phenyl}-propionic-acid
179	<p>Chiral</p>	(R)-{3-bromo-4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-acetic-acid
180	<p>Chiral</p>	(R)-3-{4-[3-(4-bromo-2-trifluoromethoxy-phenoxy)-butoxy]-2-methyl-phenyl}-propionic-acid
181	<p>Chiral</p>	(R)-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-3-methyl-phenyl}-acetic-acid
182	<p>Chiral</p>	(R)-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-acetic-acid
183	<p>Chiral</p>	(R)-3-{4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-2-trifluoromethyl-phenyl}-propionic-acid

	Structure	Name
184		(R)-14-[3-(4-chloro-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenoxy]-acetic acid
185		(R)-3-[4-[3-(4-chloro-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenyl]-propionic acid
186		(R)-3-[2-Chloro-4-[3-(4-chloro-2-phenoxy-phenoxy)-butoxy]-phenyl]-propionic acid
187		(R)-3-[4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-fluoro-phenyl]-propionic acid
188		(R)-3-[4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-ethyl-phenyl]-propionic acid
189		(R)-3-[4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-chloro-phenyl]-propionic acid

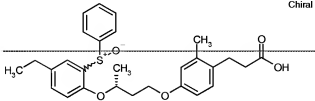
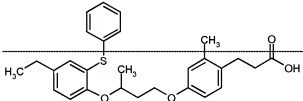
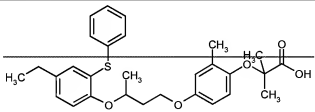
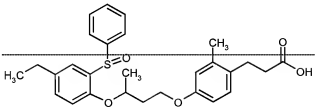
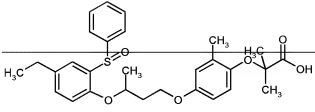
	Structure	Name
190		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-fluoro-phenyl}-propionic acid
191		(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
192		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-phenyl}-propionic acid
193		(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-pentyloxy]-2-methyl-phenyl}-propionic acid
194		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-pentyloxy]-2-methyl-phenyl}-propionic acid
195		(R)-[4-[3-(3-Benzoyl-naphthalen-2-yloxy)-butoxy]-2-methyl-phenylsulfanyl]-acetic acid

	Structure	Name
196		(R)-3-{4-[3-(2-Benzoyl-naphthalen-2-yloxy)-butoxy]-2-methyl-phenyl}-propionic-acid
197		(R)-3-{4-[3-(4-Ethyl-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic-acid
198		(R)-3-{4-[3-(4-Isopropyl-2-phenoxy-phenoxy)-butylsulfanyl]-2-methyl-phenyl}-propionic-acid
199		(R)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-propyl-phenyl}-propionic-acid
200		(R)-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-ethyl-phenylsulfanyl}-acetic-acid
201		(R)-3-{4-[3-(2-Benzoyl-4,5-dichloro-phenoxy)-butoxy]-2-methyl-phenyl}-propionic-acid

	Structure	Name
202		(R)-3-{2-Methyl-4-[3-(2-phenoxy-4-trifluoromethyl-phenoxy)-butylsulfanyl]-phenyl}-propionic acid
203		(R)-3-{2-Ethyl-4-[3-(4-ethyl-2-phenoxy-phenoxy)-butoxy]-phenyl}-propionic acid
204		(R)-3-{2-Ethyl-4-[3-(2-phenoxy-4-trifluoromethyl-phenoxy)-butoxy]-phenyl}-propionic acid
205		(R)-3-{4-[3-(2-Benzoyl-4-ethyl-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic acid
206		(R)-3-{2-Ethyl-4-[1-methyl-3-(2-phenoxy-4-trifluoromethyl-phenoxy)-propoxy]-phenyl}-propionic acid
207		(R)-3-{2-Methyl-4-[1-methyl-3-(2-phenoxy-4-trifluoromethoxy-phenoxy)-propylsulfanyl]-phenyl}-propionic acid

	Structure	Name
208	 <p>Chiral</p>	(S)-3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic-acid
209		3-{4-[3-(4-Chloro-2-phenoxy-phenoxy)-propoxy]-2-ethyl-phenyl}-propionic-acid
210	 <p>Chiral</p>	(R)-3-{4-[3-(2,4-Diphenoxy-phenoxy)-butoxy]-2-ethyl-phenyl}-propionic-acid
211	 <p>Chiral Cis - Isomer 2</p>	2-{4-[4-(4-Chloro-2-phenoxy-phenyl)-3-methyl-butoxy]-2-methyl-phenyl}-cyclopropanecarboxylic acid
212		(R,S)-2-{4-[3-(4-Ethyl-2-phenylsulfanyl-phenoxy)-butoxy]-phenoxy]-2-methyl-propionic-acid
213		2-{4-[3-(R,S-2-Benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenylsulfanyl]-2-methyl-propionic acid (enantiomer pair-1)

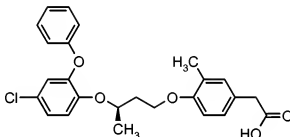
	Structure	Name
214		(R, S)-2-[4-{3-(2-cyclopropylmethyl-4-trifluoromethyl-phenoxy)-butoxy}-phenoxy]-2-methyl-propionic-acid
215		(R, S)-2-Methyl-2-[4-{3-(2-methyl-3-phenyl-7-propyl-benzofuran-6-yloxy)-butoxy}-phenoxy]-propionic-acid
216		(R, S)-2-Methyl-2-[4-{3-(4-methyl-3-phenyl-7-propyl-benzofuran-6-yloxy)-butoxy}-phenoxy]-propionic-acid
217		(R, S)-2-[4-{3-(2-cyclopropylmethyl-4-trifluoromethyl-phenoxy)-butoxy}-2-methyl-phenoxy]-2-methyl-propionic-acid
218		(R, S)-3-[4-{3-(2-cyclopropylmethyl-4-trifluoromethyl-phenoxy)-butoxy}-2-methyl-phenoxy]-2-methyl-propionic-acid

	Structure	Name
		methyl-phenyl}-propionic acid
219	 <p>Chiral</p>	3-{[R-4-{3-(R,S-2-benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
220		3-{[4-{3-(4-Ethyl-2-phenylsulfanyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid-isomer-2
221		(R,S)-2-[4-{3-(4-Ethyl-2-phenylsulfanyl-phenoxy)-butoxy]phenoxy]-2-methyl-propionic acid
222		(R,S)-3-[4-{3-(R,S-2-benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
223		(R,S)-2-[4-{3-(R,S-2-benzenesulfinyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

	Structure	Name
		propionic acid
224		(R, S)-3-{4-[3-(2-Benzenesulfonyl-4-ethyl-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid
225		3-{4-[3-(2-Benzoyl-4-(trifluoromethoxy)-phenoxy)-butoxy]-2-methyl-phenyl}-propionic acid

30. (Withdrawn). The compound of Claim 29, wherein the compound is

Chiral



or a pharmaceutically acceptable salt, solvate or hydrate thereof.

31. (Previously Presented). A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a compound of Claim 1 or a pharmaceutically acceptable salt, solvate or hydrate thereof.

32. (Canceled).
 33. (Canceled).
 34. (Canceled).
 35. (Canceled).
 36. (Canceled).

37. (Canceled).

38. (Canceled).

39. (Canceled).

40. (Canceled).

41. (Canceled).

42. (Canceled).

43. (Previously Presented). A method for lowering blood-glucose in a mammal in need thereof comprising the step of administering an effective amount of a compound of Claim 1.

44. (Canceled).

45. (Canceled).

46. (Canceled).

47. (Canceled).

48. (Canceled).

49. (Canceled)